

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025036.D
 Acq On : 25 Feb 2020 09:52
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Feb 25 11:03:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 01:09:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2324	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	9008	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	5749	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	13300	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	14719	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	16783	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	6671	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	18351	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	11161	0.407	ng/ul#	84
5) 2-Methylnaphthalene	12.14	142	7900	0.399	ng/ul	99
7) Acenaphthylene	14.04	152	13099	0.395	ng/ul	95
8) Acenaphthene	14.39	153	8912	0.400	ng/ul	98
9) Fluorene	15.38	166	10742	0.384	ng/ul	98
11) Pentachlorophenol	16.72	266	1563	0.385	ng/ul	99
12) Phenanthrene	17.11	178	18327	0.392	ng/ul	97
13) Anthracene	17.20	178	17811	0.393	ng/ul	96
15) Fluoranthene	19.13	202	23944	0.352	ng/ul	97
17) Pyrene	19.49	202	24196	0.365	ng/ul	98
18) Benzo(a)anthracene	21.24	228	25097	0.395	ng/ul	99
19) Chrysene	21.29	228	25311	0.400	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	26081	0.396	ng/ul	93
22) Benzo(k)fluoranthene	22.84	252	26600	0.408	ng/ul#	95
23) Benzo(a)pyrene	23.36	252	23886	0.398	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.65	276	30221	0.402	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	25195	0.398	ng/ul	96
26) Benzo(g,h,i)perylene	26.31	276	25582	0.403	ng/ul#	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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